Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1611txm

PASSWORD:
* * * * * * RECONNECTED TO STN INTERNATIONAL * * * * * *
SESSION RESUMED IN FILE 'REGISTRY' AT 18:20:53 ON 04 MAY 2005
FILE 'REGISTRY' ENTERED AT 18:20:53 ON 04 MAY 2005

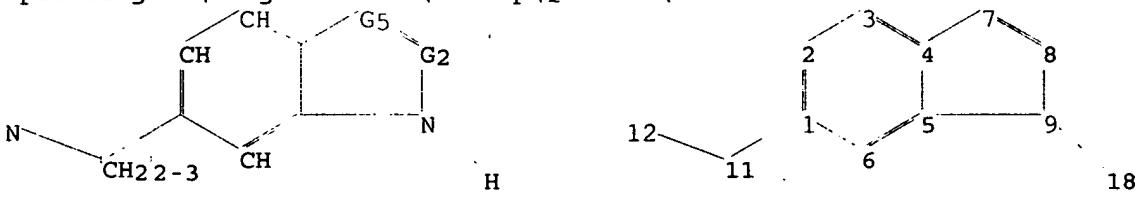
=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)es
MISSING OPERATOR

=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)/es
833648 NC4-C6/ES
53398 N2C3-C6/ES
242909 NCNC2-C6/ES
41176 N3C2-C6/ES
SYSTEM LIMITS EXCEEDED - SEARCH ENDED
The search profile you entered was too complex or gave too many answers. Simplify or subdivide the query and try again. If you have exceeded the answer limit, enter DELETE HISTORY at an arrow prompt (=>) to remove all previous answers sets and begin at L1. Use the SAVE command to store any important profiles or answer sets before using DELETE HISTORY.

=> delete history
DELETE ALL L# ITEMS? (Y)/N:.
ALL L# ITEMS DELETED

=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)/es
833648 NC4-C6/ES
53398 N2C3-C6/ES
242909 NCNC2-C6/ES
41176 N3C2-C6/ES
L1 1163389 (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES

=>
Uploading C:\Program Files\Stnexp\Queries\10784673.str



chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :

1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L2 STRUCTURE UPLOADED

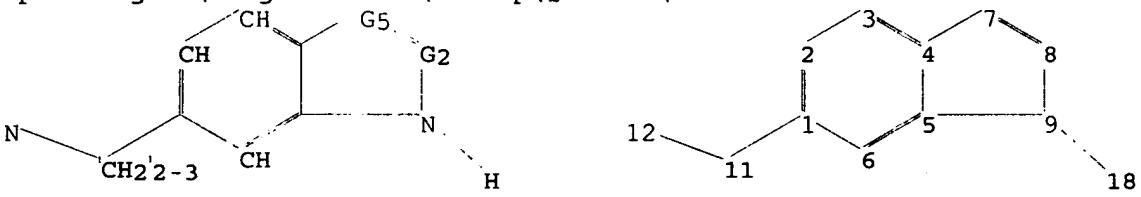
=> s l2 subset = l1 sample
SAMPLE SUBSET SEARCH INITIATED 18:23:24 FILE 'REGISTRY'
SAMPLE SUBSET SCREEN SEARCH COMPLETED - 43320 TO ITERATE

2.3% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **INCOMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 853978 TO 878822
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 0 TO 0

L3 0 SEA SUB=L1 SSS SAM L2

=>
Uploading C:\Program Files\Stnexp\Queries\10784673.str



chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :
1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12

normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L4 STRUCTURE UPLOADED

=> s l4 subset = l1 sample
SAMPLE SUBSET SEARCH INITIATED 18:24:51 FILE 'REGISTRY'
SAMPLE SUBSET SCREEN SEARCH COMPLETED - 43320 TO ITERATE

2.3% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **INCOMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 853978 TO 878822
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 0 TO 0

L5 0 SEA SUB=L1 SSS SAM L4

=>
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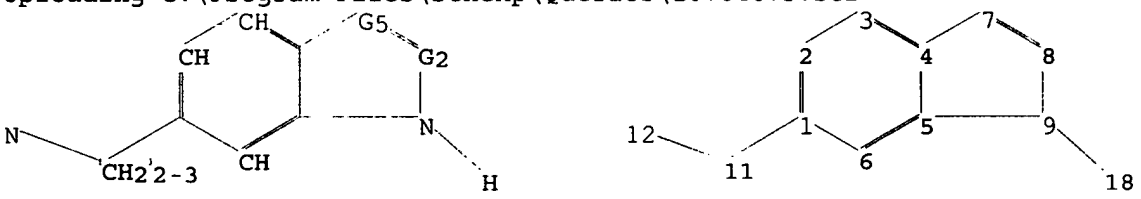


chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :
1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L6 STRUCTURE UPLOADED

=> s l6 subset = l1 sample
SAMPLE SUBSET SEARCH INITIATED 18:27:09 FILE 'REGISTRY'
SAMPLE SUBSET SCREEN SEARCH COMPLETED - 14401 TO ITERATE

6.9% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **COMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 280833 TO 295207
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 0 TO 0

L7 0 SEA SUB=L1 SSS SAM L6

=> s l6 subset = l1 full
FULL SUBSET SEARCH INITIATED 18:27:26 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 285210 TO ITERATE

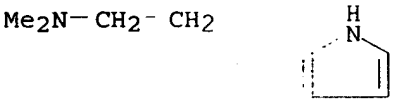
100.0% PROCESSED 285210 ITERATIONS 151 ANSWERS
SEARCH TIME: 00.00.02

L8 151 SEA SUB=L1 SSS FUL L6

=> s c12h16n2 and l8
1947 C12H16N2
L9 1 C12H16N2 AND L8

=> d ide cbib pi

L9 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN 321744-76-7 REGISTRY
ED Entered STN: 14 Feb 2001
CN 1H-Indole-6-ethanamine, N,N-dimethyl- (9CI) (CA INDEX NAME)
OTHER NAMES:
CN 6-[2-(Dimethylamino)ethyl]-1H-indole
FS 3D CONCORD
MF **C12 H16 N2**
SR CA
LC STN Files: CA, CAPLUS, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 140:303530 Preparation of indole and indazole derivatives for the treatment of migraine. Edwards, Louise; Isaac, Methvin; Maddaford, Shawn; Slassi, Abdelmalik; Xin, Tao (NPS Allelix Biopharmaceuticals, Inc., Can.). U.S. US 6716837 B1 20040406, 29 pp., Cont.-in-part of U.S. Ser. No. 354,091, abandoned. (English). CODEN: USXXAM. APPLICATION: US 2000-709579 20001113. PRIORITY: US 1999-354091 19990715.

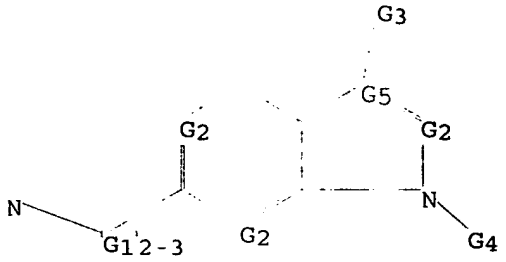
	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 6716837	B1	20040406	US 2000-709579	20001113
	US 2004167158	A1	20040826	US 2004-784673	20040224

REFERENCE 2: 134:131423 Preparation of aminoalkylindoles and analogs as 5-HT1D receptor ligands. Edwards, Louise; Isaac, Methvin; Maddaford, Shawn; Slassi, Abelmalik; Xin, Tao (NPS Allelix Corp., Can.). PCT Int. Appl. WO 2001005758 A2 20010125, 71 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 2000-CA831 20000714. PRIORITY: US 1999-354091 19990715.

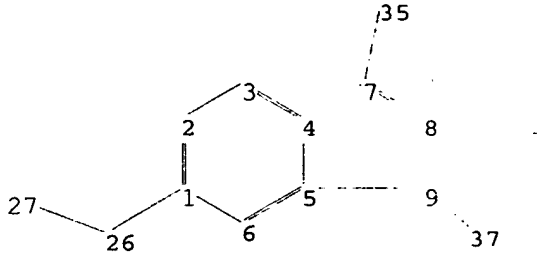
	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001005758	A2	20010125	WO 2000-CA831	20000714
	WO 2001005758	A3	20010719		
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	CA 2378047	AA	20010125	CA 2000-2378047	20000714
	EP 1196380	A2	20020417	EP 2000-945511	20000714
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
	JP 2003505369	T2	20030212	JP 2001-511419	20000714

=>
Uploading C:\Program Files\Stnexp\Queries\10784673claim7.str

Cb 9

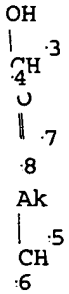


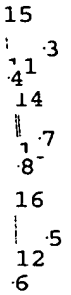
32 9



2 1
CH_2

2 1
10





chain nodes :
10 11 12 13 14 15 16 26 32 35 37
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
27
chain bonds :
1-26 7-35 9-37 11-15 12-16 13-14 26-27
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-2 1-6 1-26 2-3 3-4 4-5 4-7 5-6 5-9 7-8 7-35 8-9 9-37 11-15 12-16
13-14 26-27

G1:[*1-*2],[*3-*4],[*5-*6],[*7-*8]

G2:N,CH

G3:H,CH3,Et,n-Pr,i-Pr,n-Bu,i-Bu,s-Bu,t-Bu,[*9]

G4:i-Pr,Cy

G5:C,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 26:CLASS 27:CLASS
32:Atom 35:CLASS 37:CLASS
Generic attributes :
32:
Saturation : Saturated
Type of Ring System : Monocyclic

L10 STRUCTURE UPLOADED

=> d his

(FILE 'REGISTRY' ENTERED AT 17:37:55 ON 04 MAY 2005)
DELETE HISTORY
L1 1163389 S (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES
L2 STRUCTURE UPLOADED
L3 0 S L2 SAMPLE SUB=L1
L4 STRUCTURE UPLOADED
L5 0 S L4 SAMPLE SUB=L1
L6 STRUCTURE UPLOADED
L7 0 S L6 SAMPLE SUB=L1
L8 151 S L6 FULL SUB=L1
L9 1 S C12H16N2 AND L8
L10 STRUCTURE UPLOADED

=> s l10 subset = l8
ENTER SUBSET SEARCH SCOPE - SAMPLE, FULL, RANGE, OR (END):full
FULL SUBSET SEARCH INITIATED 18:30:13 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 151 TO ITERATE

100.0% PROCESSED 151 ITERATIONS 24 ANSWERS
SEARCH TIME: 00.00.01

L11 24 SEA SUB=L8 SSS FUL L10

=> file caold caplus;
FILE 'CAOLD' ENTERED AT 18:30:35 ON 04 MAY 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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FILE 'CAPLUS' ENTERED AT 18:30:35 ON 04 MAY 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

=> s l11

```
L12          2 L11

=> s us-6716837/pn
L13          1 US-6716837/PN

=> s l12 not l13
L14          1 L12 NOT L13

=> d cbib pi hitstr

L14  ANSWER 1 OF 1  CAPLUS  COPYRIGHT 2005 ACS on STN
2001:63967   Document No. 134:131423  Preparation of aminoalkylindoles and
     analogs as 5-HT1D receptor ligands.  Edwards, Louise; Isaac, Methvin;
     Maddaford, Shawn; Slassi, Abelmalik; Xin, Tao (NPS Allelix Corp., Can.).
     PCT Int. Appl. WO 2001005758 A2 20010125, 71 pp.  DESIGNATED STATES: W:
     AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU,
     CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN,
     IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK,
     MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM,
     TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU,
     TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR,
     GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG.
     (English).  CODEN: PIXXD2.  APPLICATION: WO 2000-CA831 20000714.
     PRIORITY: US 1999-354091 19990715.
     PATENT NO.           KIND   DATE        APPLICATION NO.         DATE
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PI   WO 2001005758         A2    20010125    WO 2000-CA831           20000714
     WO 2001005758         A3    20010719
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             CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR,
             HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT,
             LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU,
             SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN,
             YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
             DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
             CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG
     CA 2378047            AA    20010125    CA 2000-2378047         20000714
     EP 1196380            A2    20020417    EP 2000-945511          20000714
         R:  AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
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     JP 2003505369         T2    20030212    JP 2001-511419          20000714
IT   321744-77-8P 321744-79-0P 321744-81-4P
     321744-82-5P 321744-83-6P 321744-94-9P
     321744-95-0P 321744-96-1P 321745-29-3P
     321745-31-7P 321745-32-8P 321745-33-9P
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     321745-39-5P 321745-41-9P 321745-45-3P
     321745-47-5P 321745-49-7P
     RL: BAC (Biological activity or effector, except adverse); BSU (Biological
     study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
     BIOL (Biological study); PREP (Preparation); USES (Uses)
        (preparation of aminoalkylindoles and analogs as 5-HT1D receptor ligands)
RN   321744-77-8  CAPLUS
CN   1H-Indole-6-ethanamine, N,N-dimethyl-1-(1-methylethyl)- (9CI)  (CA INDEX
     NAME)
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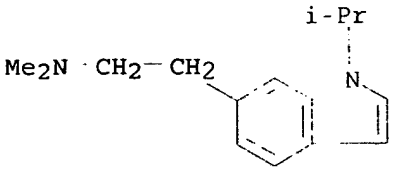
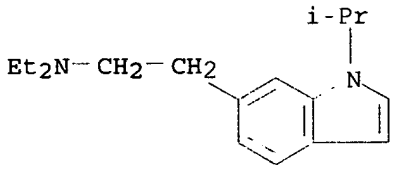
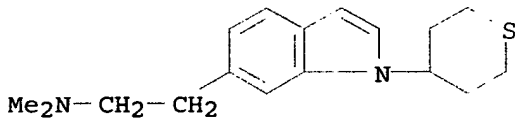


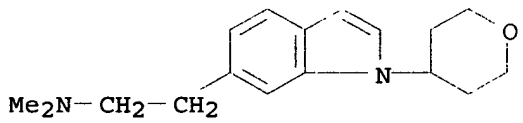
RN 321744-79-0 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



RN 321744-81-4 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(tetrahydro-2H-thiopyran-4-yl)- (9CI) (CA INDEX NAME)



RN 321744-82-5 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(tetrahydro-2H-pyran-4-yl)- (9CI) (CA INDEX NAME)



RN 321744-83-6 CAPLUS
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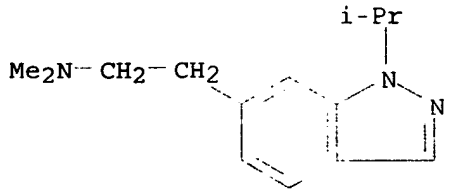
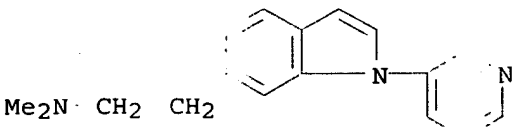


RN 321744-94-9 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(3-pyridinyl)- (9CI) (CA INDEX

NAME)



RN 321744-95-0 CAPLUS
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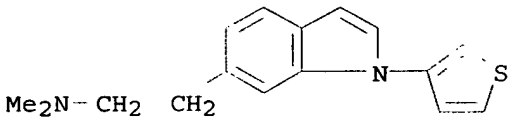
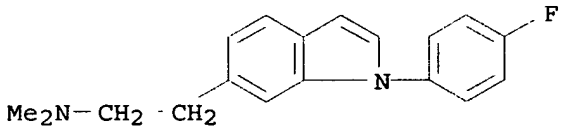
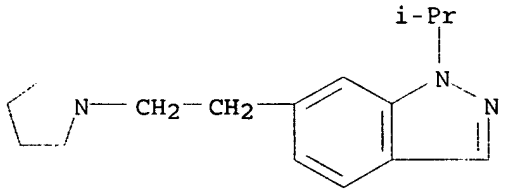


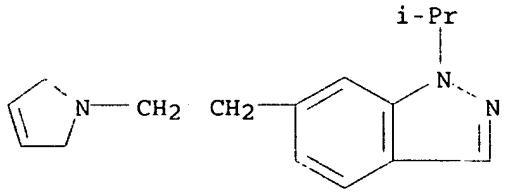
RN 321744-96-1 CAPLUS
CN 1H-Indole-6-ethanamine, 1-(4-fluorophenyl)-N,N-dimethyl- (9CI) (CA INDEX NAME)



RN 321745-29-3 CAPLUS
CN 1H-Indazole, 1-(1-methylethyl)-6-[2-(1-pyrrolidinyl)ethyl]- (9CI) (CA INDEX NAME)

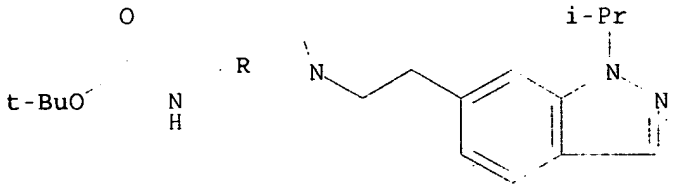


RN 321745-31-7 CAPLUS
CN 1H-Indazole, 6-[2-(2,5-dihydro-1H-pyrrol-1-yl)ethyl]-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



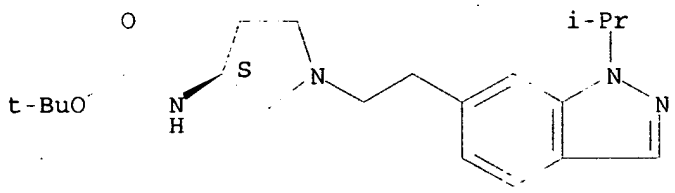
RN 321745-32-8 CAPLUS
CN Carbamic acid, [(3R)-1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-3-pyrrolidinyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.



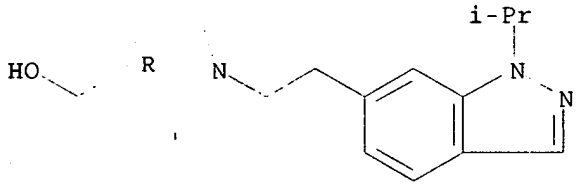
RN 321745-33-9 CAPLUS
CN Carbamic acid, [(3S)-1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-3-pyrrolidinyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.



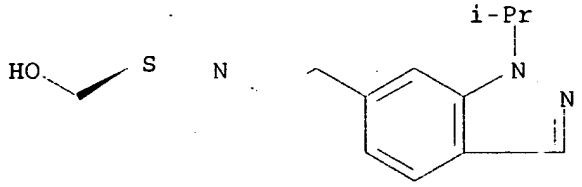
RN 321745-34-0 CAPLUS
CN 3-Pyrrolidinemethanol, 1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-, (3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

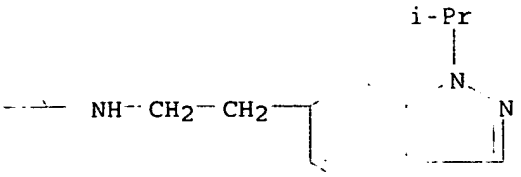


RN 321745-35-1 CAPLUS
CN 3-Pyrrolidinemethanol, 1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-, (3S)- (9CI) (CA INDEX NAME)

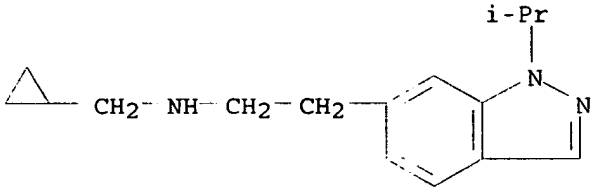
Absolute stereochemistry.



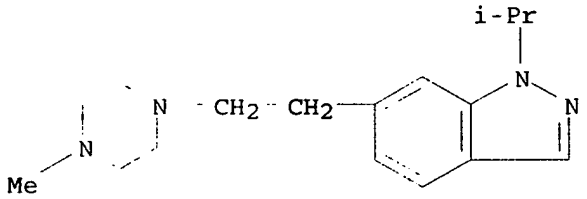
RN 321745-37-3 CAPLUS
CN 1H-Indazole-6-ethanamine, N-cyclopropyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



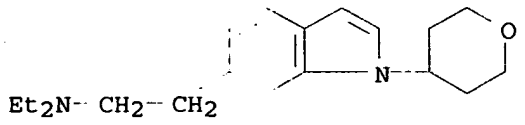
RN 321745-39-5 CAPLUS
CN 1H-Indazole-6-ethanamine, N-(cyclopropylmethyl)-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



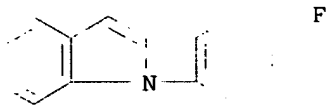
RN 321745-41-9 CAPLUS
CN 1H-Indazole, 1-(1-methylethyl)-6-[2-(4-methyl-1-piperazinyl)ethyl]- (9CI) (CA INDEX NAME)



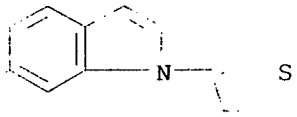
RN 321745-45-3 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(tetrahydro-2H-pyran-4-yl)- (9CI) (CA INDEX NAME)



RN 321745-47-5 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(4-fluorophenyl)- (9CI) (CA INDEX NAME)



RN 321745-49-7 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(3-thienyl)- (9CI) (CA INDEX NAME)



=> d his

(FILE 'REGISTRY' ENTERED AT 17:37:55 ON 04 MAY 2005)
DELETE HISTORY
L1 1163389 S (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES
L2 STRUCTURE UPLOADED
L3 0 S L2 SAMPLE SUB=L1
L4 STRUCTURE UPLOADED
L5 0 S L4 SAMPLE SUB=L1
L6 STRUCTURE UPLOADED
L7 0 S L6 SAMPLE SUB=L1
L8 151 S L6 FULL SUB=L1
L9 1 S C12H16N2 AND L8
L10 STRUCTURE UPLOADED
L11 24 S L10 SUB=L8 FULL

FILE 'CAOLD, CAPLUS' ENTERED AT 18:30:35 ON 04 MAY 2005
L12 2 S L11
L13 1 S US-6716837/PN
L14 1 S L12 NOT L13

=> s l8
L15 58 L8

=> s l15 not l12
L16 56 L15 NOT L12

=> sort py l16
SORT ENTIRE ANSWER SET? (Y)/N:.
2 ANSWERS DID NOT HAVE 'PY' SORT FIELD
PROCESSING COMPLETED FOR L16
L17 56 SORT L16 PY

=> file reg
FILE 'REGISTRY' ENTERED AT 18:33:27 ON 04 MAY 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file

provided by InfoChem.

STRUCTURE FILE UPDATES: 3 MAY 2005 HIGHEST RN 849720-40-7
DICTIONARY FILE UPDATES: 3 MAY 2005 HIGHEST RN 849720-40-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

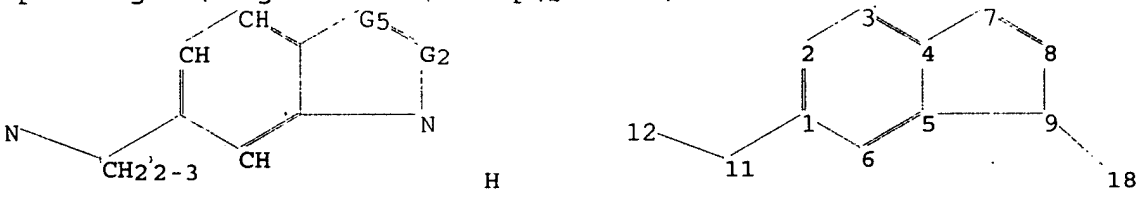
Please note that search-term pricing does apply when conducting SmartSELECT searches.

****
* *
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
* *
****

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Program Files\Stnexp\Queries\10784673.str



chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :
1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L18 STRUCTURE UPLOADED

=> s l18 subset = l8 full
FULL SUBSET SEARCH INITIATED 18:33:55 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 151 TO ITERATE

100.0% PROCESSED 151 ITERATIONS 113 ANSWERS
SEARCH TIME: 00.00.01

L19 113 SEA SUB=L8 SSS FUL L18

=> s l8 not l19
L20 38 L8 NOT L19

=> file caold caplus; s l20
FILE 'CAOLD' ENTERED AT 18:34:31 ON 04 MAY 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
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COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'CAPLUS' ENTERED AT 18:34:31 ON 04 MAY 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

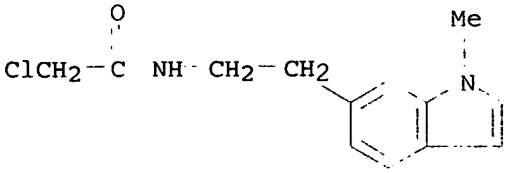
L21 8 L20

=> s l21 not l12
L22 6 L21 NOT L12

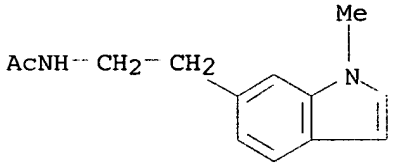
=> sort py l22
SORT ENTIRE ANSWER SET? (Y)/N:.
PROCESSING COMPLETED FOR L22
L23 6 SORT L22 PY

=> d 1-6 cbib pi fhitstr

L23 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1976:16500 Document No. 84:16500 Photocyclization of N-chloroacetyl derivatives of indolylethylamines to azepinoindoles and azocinoindoles. Correlation of the reactivity of indole radicals with frontier electron densities calculated by unrestricted Hartree-Fock MO. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Tokyo, Japan). Tetrahedron Letters (39), 3399-402 (English) 1975. CODEN: TELEAY. ISSN: 0040-4039.
IT **57964-71-3**
RL: RCT (Reactant); RACT (Reactant or reagent)
(photochem. cyclization of, MO calcns. for)
RN 57964-71-3 CAPLUS
CN Acetamide, 2-chloro-N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



L23 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1980:549357 Document No. 93:149357 Photochemical synthesis of azepinoindoles and azocinoindoles from N-chloroacetylindolylethylamines, and a mechanistic study based on the correlation between quantum yields and calculated frontier electron densities of indole radicals. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Sankyo Co., Tokyo, 140, Japan). Chemical & Pharmaceutical Bulletin, 28(3), 900-9 (English) 1980. CODEN: CPBTAL. ISSN: 0009-2363.
IT **74631-90-6P**
RL: FORM (Formation, nonpreparative); PREP (Preparation)
(formation of, by photolysis of chloroacetylindolylethylamine)
RN 74631-90-6 CAPLUS
CN Acetamide, N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



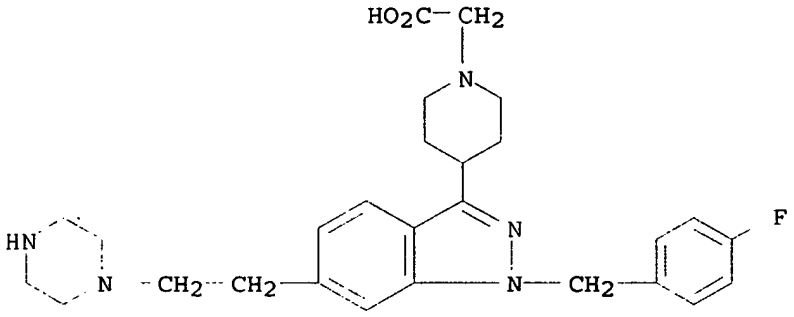
L23 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1998:42394 Document No. 128:102084 Preparation of 4-heterocyclyl-1-piperidineacetates as glycoprotein IIb/IIIa receptor antagonists. Allen, David George; Eldred, Colin David; Judkins, Brian David; Mitchell, William Leonard; Scopes, David Ian Carter (Glaxo Group Ltd., UK; Allen, David George; Eldred, Colin David; Judkins, Brian David; Mitchell, William Leonard; Scopes, David Ian Carter). PCT Int. Appl. WO 9749698 A1 19971231, 84 pp. DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 1997-EP3194 19970619. PRIORITY: GB 1996-13017 19960621; GB 1996-13095 19960621; GB 1996-13018 19960621; GB 1996-13026 19960621.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 9749698	A1	19971231	WO 1997-EP3194	19970619

W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG

AU 9732610	A1	19980114	AU 1997-32610	19970619
ZA 9705431	A	19981221	ZA 1997-5431	19970619
CN 1222153	A	19990707	CN 1997-195652	19970619

IT **201482-63-5P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of 4-heterocyclyl-1-piperidineacetates as glycoprotein IIb/IIIa receptor antagonists)
RN 201482-63-5 CAPLUS
CN 1-Piperidineacetic acid, 4-[1-[(4-fluorophenyl)methyl]-6-[2-(1-piperazinyl)ethyl]-1H-indazol-3-yl]-, tetrahydrochloride (9CI) (CA INDEX NAME)



●4 HCl

L23 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1999:311189 Document No. 130:338114 Indazole bioisostere replacement of catechol in therapeutically active compounds. Marfat, Anthony (Pfizer Products Inc., USA). PCT Int. Appl. WO 9923077 A1 19990514, 231 pp.
DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.
APPLICATION: WO 1998-IB1710 19981026. PRIORITY: US 1997-64229 19971104; US 1997-64187 19971104; US 1997-64024 19971104; US 1997-64228 19971104; US 1997-64198 19971104.

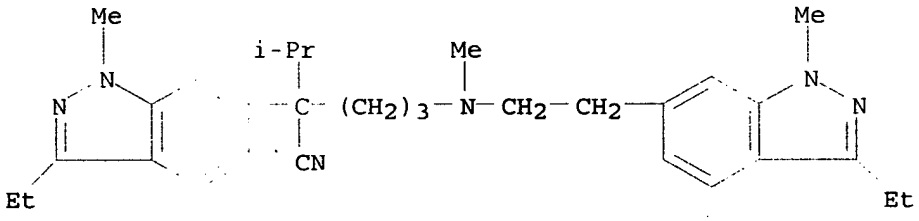
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 9923077	A1	19990514	WO 1998-IB1710	19981026
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
CA 2309150	AA	19990514	CA 1998-2309150	19981026

AU 9894552	A1	19990524	AU 1998-94552	19981026
AU 754734	B2	20021121		
TR 200001234	T2	20000821	TR 2000-200001234	19981026
EP 1028946	A1	20000823	EP 1998-947732	19981026
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE, SI, LT, LV, FI, RO				
BR 9813926	A	20000919	BR 1998-13926	19981026
JP 2001521926	T2	20011113	JP 2000-518952	19981026
NZ 503918	A	20020328	NZ 1998-503918	19981026
ZA 9810041	A	20000503	ZA 1998-10041	19981103
AP 910	A	20001205	AP 1998-1375	19981104
W: BW, GM, GH, KE, MW, SD, UG, ZM, ZW				
US 6391872	B1	20020521	US 1999-381425	19990920
US 6329412	B1	20011211	US 2000-535359	20000324
NO 2000002129	A	20000703	NO 2000-2129	20000426
HR 2000000253	A1	20010630	HR 2000-253	20000428
BG 104450	A	20001229	BG 2000-104450	20000517

IT **224188-05-0P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(indazole bioisostere replacement of catechol in therapeutically active compds.)

RN 224188-05-0 CAPLUS

CN 1H-Indazole-6-acetonitrile, 3-ethyl-α-[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-1-methyl-α-(1-methylethyl)-(9CI) (CA INDEX NAME)



L23 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

2000:68450 Document No. 132:107953 Preparation of heterocyclic compounds as antagonists of gonadotropin releasing hormone. Goulet, Mark; Ashton, Wallace T.; Chu, Lin; Fisher, Michael H.; Lin, Peter; Ponpipom, Mitree M.; Wyvratt, Matthew J.; Girotra, Narindar N.; Young, Jonathan (Merck and Co., Inc., USA). PCT Int. Appl. WO 2000004013 A1 20000127, 138 pp. DESIGNATED STATES: W: AE, AL, AM, AU, AZ, BA, BB, BG, BR, BY, CA, CN, CU, CZ, EE, GD, GE, HR, HU, ID, IL, IN, IS, JP, KG, KR, KZ, LC, LK, LR, LT, LV, MD, MG, MK, MN, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TJ, TM, TR, TT, UA, US, UZ, VN, YU, ZA, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.
APPLICATION: WO 1999-US15581 19990709. PRIORITY: US 1998-115497 19980714.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2000004013	A1	20000127	WO 1999-US15581	19990709
W: AE, AL, AM, AU, AZ, BA, BB, BG, BR, BY, CA, CN, CU, CZ, EE, GD, GE, HR, HU, ID, IL, IN, IS, JP, KG, KR, KZ, LC, LK, LR, LT, LV, MD, MG, MK, MN, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TJ, TM, TR, TT, UA, US, UZ, VN, YU, ZA, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				

RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

US 6200957	B1	20010313	US 1998-115497	19980714
CA 2337407	AA	20000127	CA 1999-2337407	19990709
AU 9949816	A1	20000207	AU 1999-49816	19990709
AU 760030	B2	20030508		
EP 1095038	A1	20010502	EP 1999-933850	19990709
EP 1095038	B1	20030903		

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO

JP 2002520409	T2	20020709	JP 2000-560119	19990709
AT 248832	E	20030915	AT 1999-933850	19990709

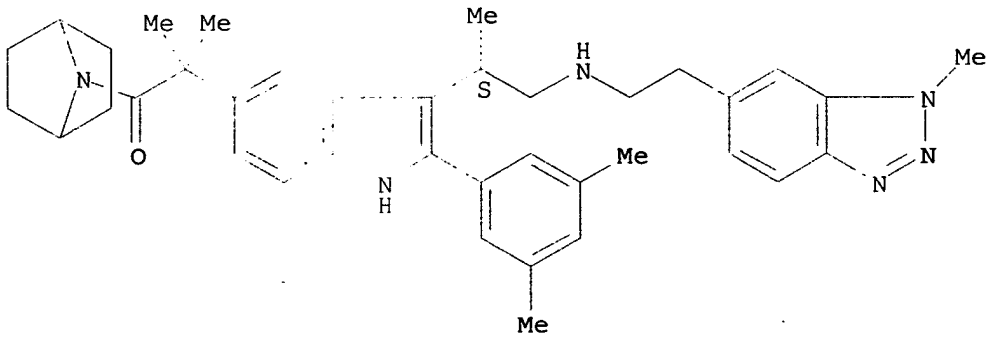
IT **255863-35-5P**

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of heterocyclic compds. as antagonists of gonadotropin releasing hormone)

RN 255863-35-5 CAPLUS

CN 7-Azabicyclo[2.2.1]heptane, 7-[2-[2-(3,5-dimethylphenyl)-3-[(1S)-1-methyl-2-[[2-(1-methyl-1H-benzotriazol-6-yl)ethyl]amino]ethyl]-1H-indol-5-yl]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L23 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

2001:178434 Document No. 134:222629 Preparation of indoles as antagonists of gonadotropin releasing hormone. Goulet, Mark; Ashton, Wallace T.; Chu, Lin; Fisher, Michael H.; Lin, Peter; Ponpipom, Mitree M.; Wyvratt, Matthew J.; Girotra, Narindar N.; Young, Jonathan (Merck & Co., Inc., USA). U.S. US 6200957 B1 20010313, 53 pp., Cont.-in-part of U.S. 5,780,437. (English). CODEN: USXXAM. APPLICATION: US 1998-115497 19980714. PRIORITY: US 1995-PV8633 19951214; US 1996-760816 19961205.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI US 6200957	B1	20010313	US 1998-115497	19980714
US 5780437	A	19980714	US 1996-760816	19961205
JP 2001106685	A2	20010417	JP 2000-257791	19961210
ZA 9610536	A	19970814	ZA 1996-10536	19961213
CA 2337407	AA	20000127	CA 1999-2337407	19990709
WO 2000004013	A1	20000127	WO 1999-US15581	19990709

W: AE, AL, AM, AU, AZ, BA, BB, BG, BR, BY, CA, CN, CU, CZ, EE, GD, GE, HR, HU, ID, IL, IN, IS, JP, KG, KR, KZ, LC, LK, LR, LT, LV, MD, MG, MK, MN, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TJ, TM, TR,

TT, UA, US, UZ, VN, YU, ZA, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

AU 9949816	A1	20000207	AU 1999-49816	19990709
AU 760030	B2	20030508		
EP 1095038	A1	20010502	EP 1999-933850	19990709
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JP 2002520409	T2	20020709	JP 2000-560119	19990709
AT 248832	E	20030915	AT 1999-933850	19990709
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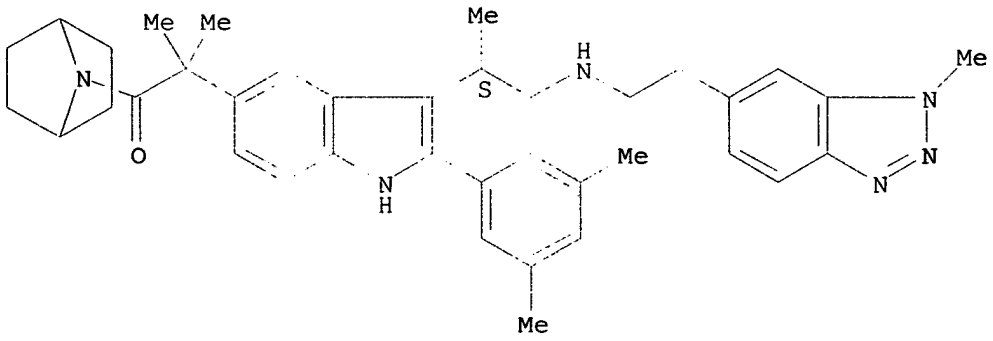
IT **255863-35-5P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of indoles as antagonists of gonadotropin releasing hormone)
RN 255863-35-5 CAPLUS
CN 7-Azabicyclo[2.2.1]heptane, 7-[2-[2-(3,5-dimethylphenyl)-3-[(1S)-1-methyl-2-[[2-(1-methyl-1H-benzotriazol-6-yl)ethyl]amino]ethyl]-1H-indol-5-yl]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



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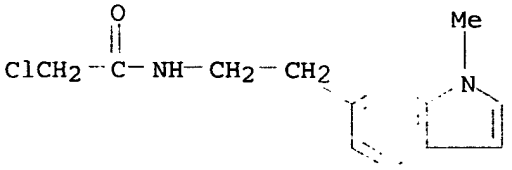
L23 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1976:16500 Document No. 84:16500 Photocyclization of N-chloroacetyl derivatives of indolylethylamines to azepinoindoles and azocinoindoles. Correlation of the reactivity of indole radicals with frontier electron densities calculated by unrestricted Hartree-Fock MO. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Tokyo, Japan). Tetrahedron Letters (39), 3399-402 (English) 1975. CODEN: TELEAY. ISSN: 0040-4039.
IT **57964-71-3**
RL: RCT (Reactant); RACT (Reactant or reagent)
(photochem. cyclization of, MO calcns. for)
RN 57964-71-3 CAPLUS
CN Acetamide, 2-chloro-N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



L23 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN
1999:311189 Document No. 130:338114 Indazole bioisostere replacement of catechol in therapeutically active compounds. Marfat, Anthony (Pfizer Products Inc., USA). PCT Int. Appl. WO 9923077 A1 19990514, 231 pp. DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 1998-IB1710 19981026. PRIORITY: US 1997-64229 19971104; US 1997-64187 19971104; US 1997-64024 19971104; US 1997-64228 19971104; US 1997-64198 19971104.

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9923077	A1	19990514	WO 1998-IB1710	19981026
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	RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	CA 2309150	AA	19990514	CA 1998-2309150	19981026
	AU 9894552	A1	19990524	AU 1998-94552	19981026
	AU 754734	B2	20021121		
	TR 200001234	T2	20000821	TR 2000-200001234	19981026
	EP 1028946	A1	20000823	EP 1998-947732	19981026
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE, SI, LT, LV, FI, RO				
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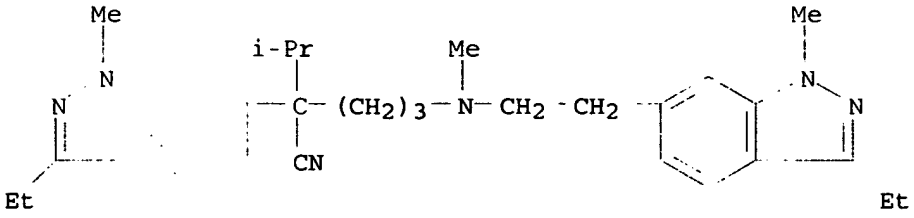
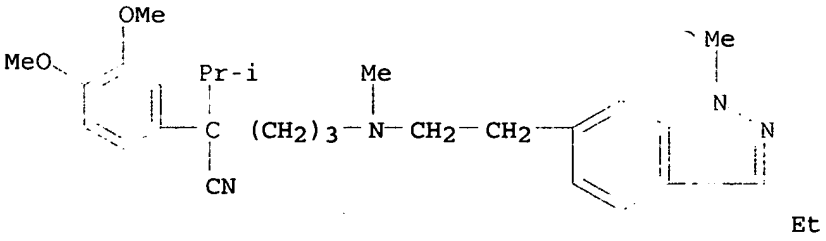
	JP 2001521926	T2	20011113	JP 2000-518952	19981026
	NZ 503918	A	20020328	NZ 1998-503918	19981026
	ZA 9810041	A	20000503	ZA 1998-10041	19981103
	AP 910	A	20001205	AP 1998-1375	19981104
	W: BW, GM, GH, KE, MW, SD, UG, ZM, ZW				
	US 6391872	B1	20020521	US 1999-381425	19990920
	US 6329412	B1	20011211	US 2000-535359	20000324
	NO 2000002129	A	20000703	NO 2000-2129	20000426
	HR 2000000253	A1	20010630	HR 2000-253	20000428
	BG 104450	A	20001229	BG 2000-104450	20000517

IT **224188-05-0P 224188-16-3P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(indazole bioisostere replacement of catechol in therapeutically active compds.)
RN 224188-05-0 CAPLUS

CN 1H-Indazole-6-acetonitrile, 3-ethyl-α-[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-1-methyl-α-(1-methylethyl)- (9CI) (CA INDEX NAME)



RN 224188-16-3 CAPLUS
CN Benzeneacetonitrile, α-[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-3,4-dimethoxy-α-(1-methylethyl)- (9CI) (CA INDEX NAME)



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L2	13	1 and (indol$ OR indazol$) and ("5-ht-1d" OR 5ht".sub.1d")	US-PGPUB; USPAT	OR	ON	2005/05/04 16:44
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